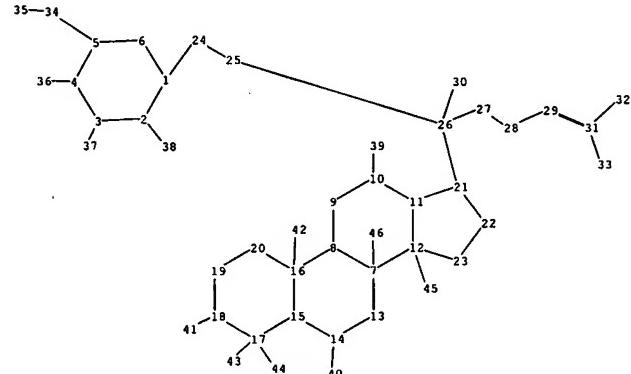
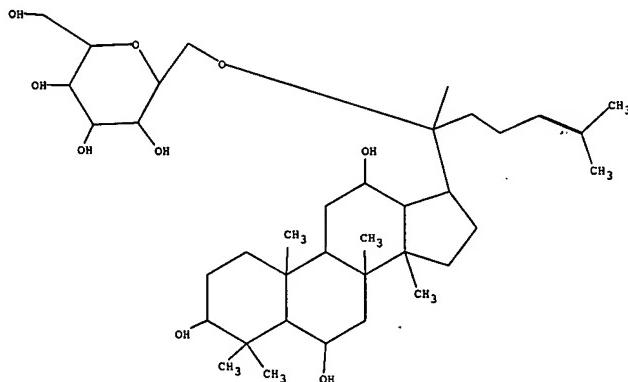


(FILE 'HOME' ENTERED AT 12:32:45 ON 31 DEC 2007)

FILE 'REGISTRY' ENTERED AT 12:33:01 ON 31 DEC 2007

L1 STRUCTURE UPLOADED
L2 0 S L1 SSS SAM
L3 0 S L1 SSS FULL



chain nodes :

24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43
44 45 46

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22
23

chain bonds :

1-24 2-38 3-37 4-36 5-34 7-46 10-39 12-45 14-40 16-42 17-43 17-44
18-41 21-26 24-25 25-26 26-27 26-30 27-28 28-29 29-31 31-32 31-33
34-35

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 7-13 8-9 8-16 9-10 10-11 11-12
11-21 12-23 13-14 14-15 15-16 15-17 16-20 17-18 18-19 19-20 21-22
22-23

exact/norm bonds :

1-2 1-6 2-3 2-38 3-4 3-37 4-5 4-36 5-6 7-8 7-12 7-13 8-9 8-16
9-10 10-11 10-39 11-12 11-21 12-23 13-14 14-15 14-40 15-16 15-17
16-20 17-18 18-19 18-41 19-20 21-22 22-23 24-25 25-26 34-35

exact bonds :

1-24 5-34 7-46 12-45 16-42 17-43 17-44 21-26 26-27 26-30 27-28 28-29
29-31 31-32 31-33

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS25:CLASS26:CLASS27:CLASS
28:CLASS29:CLASS30:CLASS31:CLASS32:CLASS33:CLASS34:CLASS35:CLASS
36:CLASS37:CLASS38:CLASS39:CLASS40:CLASS41:CLASS42:CLASS43:CLASS
44:CLASS45:CLASS46:CLASS